



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 14, 2020 – 05:46 am BST

PDB ID : 1PXH
Title : Crystal structure of protein tyrosine phosphatase 1B with potent and selective bidentate inhibitor compound 2
Authors : Sun, J.P.; Fedorov, A.; Lee, S.Y.; Guo, X.L.; Shen, K.; Lawrence, D.S.; Almo, S.C.; Zhang, Z.Y.
Deposited on : 2003-07-04
Resolution : 2.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

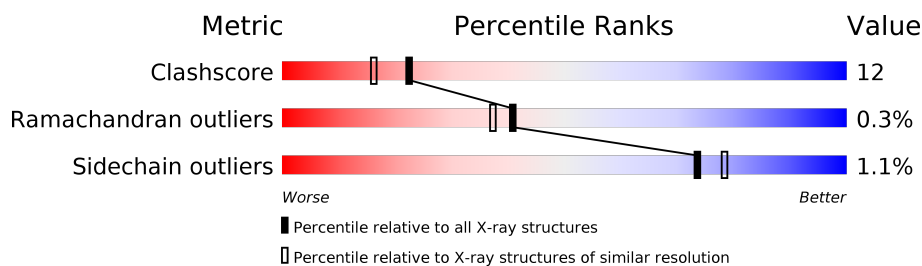
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	321	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2593 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

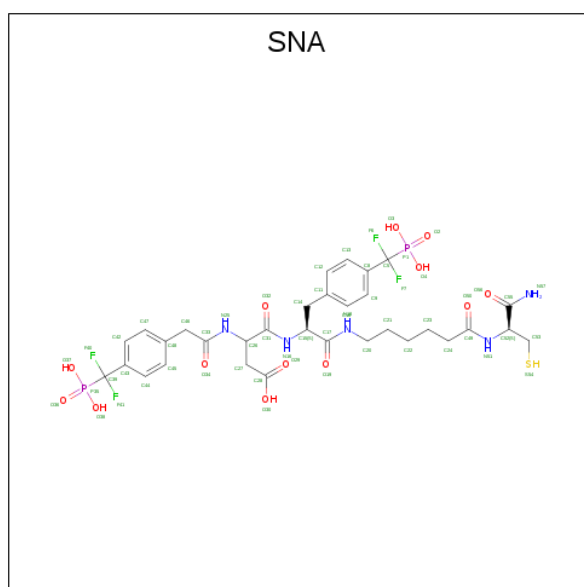
- Molecule 1 is a protein called Protein-tyrosine phosphatase, non-receptor type 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2426	1535	418	457	16			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

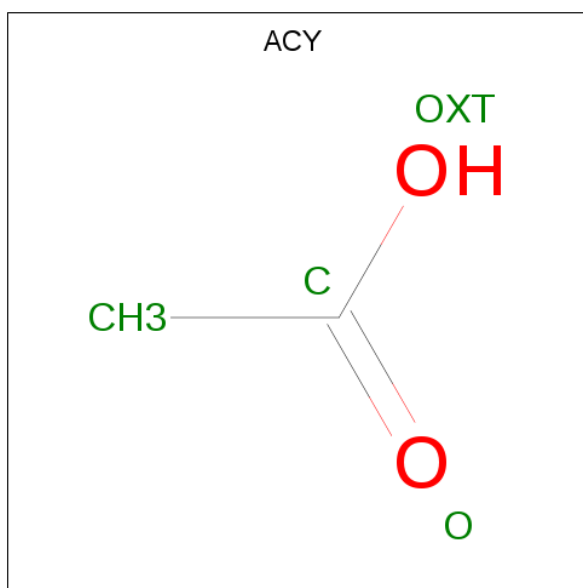
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is N-{1-[5-(1-CARBAMOYL-2-MERCAPTO-ETHYLCARBAMOYL)-PENTYL CARBAMOYL]-2-[4-(DIFLUORO-PHOSPHONO-METHYL)-PHENYL]-ETHYL}-3-{2-[4-(DIFLUORO-PHOSPHONO-METHYL)-PHENYL]-ACETYLAMINO}-SUCCINAMIC ACID (three-letter code: SNA) (formula: C₃₂H₄₁F₄N₅O₁₃P₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	P	0	0
			48	28	4	3	11	2		

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	109	Total	O	0	0
			109	109		

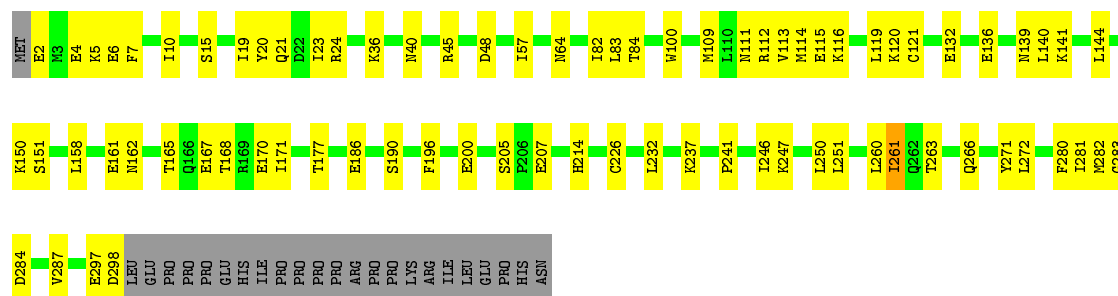
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: Protein-tyrosine phosphatase, non-receptor type 1

Chain A: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.84Å 85.68Å 88.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.15)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.207	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2593	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, MG, SNA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.32	0/2481	0.55	0/3344

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2426	0	2381	57	0
2	A	2	0	0	0	0
3	A	48	0	26	1	0
4	A	8	0	6	0	0
5	A	109	0	0	2	0
All	All	2593	0	2413	57	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (57) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:LYS:O	1:A:251:LEU:HD23	1.85	0.77
1:A:40:ASN:HD21	1:A:64:ASN:HD21	1.34	0.74
1:A:40:ASN:ND2	1:A:64:ASN:HD21	1.89	0.69
1:A:2:GLU:HG2	1:A:4:GLU:H	1.58	0.68
1:A:151:SER:N	1:A:298:ASP:HB2	2.08	0.67
1:A:100:TRP:CH2	1:A:167:GLU:HG3	2.30	0.66
1:A:115:GLU:O	1:A:116:LYS:HG2	1.98	0.64
1:A:263:THR:OG1	1:A:266:GLN:HG3	2.01	0.60
1:A:150:LYS:HA	1:A:298:ASP:HB3	1.84	0.60
1:A:100:TRP:CZ2	1:A:167:GLU:HG3	2.36	0.59
1:A:111:ASN:HB2	1:A:115:GLU:OE2	2.02	0.59
1:A:116:LYS:HD3	1:A:120:LYS:NZ	2.17	0.59
1:A:140:LEU:HD23	1:A:162:ASN:HA	1.85	0.59
1:A:237:LYS:HG3	5:A:572:HOH:O	2.06	0.56
1:A:246:ILE:HG22	5:A:537:HOH:O	2.05	0.56
1:A:232:LEU:HB3	1:A:281:ILE:HD11	1.87	0.55
1:A:2:GLU:OE1	1:A:4:GLU:HB3	2.08	0.54
1:A:165:THR:OG1	1:A:167:GLU:HG2	2.09	0.53
1:A:196:PHE:O	1:A:200:GLU:HG3	2.09	0.53
1:A:205:SER:OG	1:A:207:GLU:HG2	2.09	0.52
1:A:241:PRO:HG3	1:A:282:MET:CE	2.39	0.52
1:A:150:LYS:HD2	1:A:297:GLU:HB3	1.91	0.52
1:A:161:GLU:HB2	1:A:168:THR:HG22	1.92	0.51
1:A:113:VAL:HG13	1:A:121:CYS:O	2.11	0.51
1:A:170:GLU:C	1:A:171:ILE:HD12	2.32	0.50
1:A:112:ARG:HD2	1:A:177:THR:O	2.12	0.50
1:A:36:LYS:HA	1:A:36:LYS:HE2	1.93	0.50
1:A:15:SER:O	1:A:19:ILE:HD13	2.12	0.49
1:A:114:MET:HG2	1:A:119:LEU:CD2	2.42	0.49
1:A:115:GLU:C	1:A:116:LYS:HG2	2.33	0.48
1:A:186:GLU:HG2	1:A:190:SER:OG	2.14	0.47
1:A:100:TRP:HH2	1:A:167:GLU:HG3	1.80	0.47
1:A:20:TYR:CE2	1:A:24:ARG:HD2	2.50	0.46
1:A:132:GLU:HG3	1:A:141:LYS:HE3	1.98	0.46
1:A:171:ILE:HD12	1:A:171:ILE:N	2.31	0.46
1:A:151:SER:H	1:A:298:ASP:H	1.62	0.46
1:A:23:ILE:HD13	1:A:250:LEU:HD23	1.98	0.45
1:A:109:MET:HG3	1:A:214:HIS:CE1	2.51	0.45
1:A:260:LEU:O	1:A:261:ILE:HB	2.16	0.45
1:A:283:GLY:O	1:A:284:ASP:HB3	2.17	0.45
1:A:6:GLU:O	1:A:10:ILE:HG13	2.17	0.44
1:A:150:LYS:C	1:A:298:ASP:HB2	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:ILE:HD11	1:A:82:ILE:HD12	1.99	0.43
1:A:116:LYS:HD3	1:A:120:LYS:HZ1	1.81	0.43
1:A:280:PHE:HB2	1:A:287:VAL:CG2	2.48	0.43
1:A:83:LEU:HD11	1:A:226:CYS:SG	2.59	0.43
1:A:84:THR:O	1:A:214:HIS:HB2	2.19	0.42
1:A:144:LEU:HA	1:A:158:LEU:HD23	2.01	0.42
1:A:48:ASP:CG	3:A:401:SNA:H271	2.39	0.42
1:A:246:ILE:HG21	1:A:271:TYR:CZ	2.54	0.42
1:A:45:ARG:NH2	1:A:121:CYS:HA	2.35	0.41
1:A:57:ILE:HD11	1:A:82:ILE:CD1	2.50	0.41
1:A:7:PHE:CZ	1:A:272:LEU:HD21	2.56	0.41
1:A:19:ILE:N	1:A:19:ILE:HD12	2.36	0.41
1:A:246:ILE:CG2	1:A:271:TYR:CZ	3.04	0.41
1:A:21:GLN:HE22	1:A:24:ARG:HE	1.69	0.41
1:A:241:PRO:HG3	1:A:282:MET:HE1	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	295/321 (92%)	283 (96%)	11 (4%)	1 (0%)	41 37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	261	ILE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	270 / 294 (92%)	267 (99%)	3 (1%)	73	78

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	136	GLU
1	A	139	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	40	ASN
1	A	139	ASN
1	A	288	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SNA	A	401	-	40,49,58	6.62	24 (60%)	58,72,84	2.74	28 (48%)
4	ACY	A	402	-	1,3,3	5.96	1 (100%)	0,3,3	0.00	-
4	ACY	A	403	-	1,3,3	5.64	1 (100%)	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SNA	A	401	-	-	7/52/64/77	0/2/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	SNA	C44-C43	14.94	1.63	1.39
3	A	401	SNA	C39-C43	14.60	1.63	1.50
3	A	401	SNA	C9-C8	13.83	1.61	1.39
3	A	401	SNA	C42-C43	12.34	1.59	1.39
3	A	401	SNA	C13-C8	11.82	1.58	1.39
3	A	401	SNA	C46-C48	-9.39	1.36	1.51
3	A	401	SNA	P35-O36	8.76	1.64	1.50
3	A	401	SNA	C45-C48	8.70	1.57	1.38
3	A	401	SNA	P35-O38	8.22	1.69	1.54
3	A	401	SNA	C47-C48	7.70	1.55	1.38
3	A	401	SNA	C47-C42	7.56	1.52	1.38
3	A	401	SNA	C45-C44	7.55	1.52	1.38
3	A	401	SNA	O19-C17	7.16	1.37	1.23
3	A	401	SNA	C10-C11	7.12	1.54	1.38
3	A	401	SNA	C14-C11	-7.03	1.34	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	SNA	C5-C8	6.87	1.56	1.50
3	A	401	SNA	C13-C12	6.07	1.49	1.38
3	A	401	SNA	C46-C33	5.97	1.65	1.51
4	A	402	ACY	CH3-C	5.96	1.56	1.48
4	A	403	ACY	CH3-C	5.64	1.55	1.48
3	A	401	SNA	C20-N18	4.50	1.56	1.46
3	A	401	SNA	C12-C11	3.84	1.47	1.38
3	A	401	SNA	C21-C20	3.41	1.65	1.51
3	A	401	SNA	C10-C9	2.78	1.43	1.38
3	A	401	SNA	C17-N18	2.72	1.39	1.33
3	A	401	SNA	P1-O2	2.45	1.54	1.50

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	SNA	C27-C26-N25	7.46	126.51	110.79
3	A	401	SNA	O19-C17-C15	-6.35	107.09	120.45
3	A	401	SNA	C46-C33-N25	-5.41	107.17	115.88
3	A	401	SNA	F6-C5-C8	-4.60	104.56	110.49
3	A	401	SNA	C14-C11-C10	-4.49	111.99	120.91
3	A	401	SNA	O19-C17-N18	-4.42	113.51	122.99
3	A	401	SNA	C27-C26-C31	4.36	121.57	110.25
3	A	401	SNA	O32-C31-N16	3.95	130.24	122.93
3	A	401	SNA	O34-C33-C46	-3.78	113.42	122.03
3	A	401	SNA	C26-C31-N16	-3.68	108.62	116.70
3	A	401	SNA	C12-C11-C10	3.66	123.92	118.17
3	A	401	SNA	C21-C20-N18	-3.46	102.31	112.21
3	A	401	SNA	C14-C15-C17	3.44	119.19	110.25
3	A	401	SNA	F41-C39-C43	-3.36	106.16	110.49
3	A	401	SNA	C11-C14-C15	3.15	122.07	113.39
3	A	401	SNA	C12-C13-C8	-3.13	117.03	121.22
3	A	401	SNA	C17-C15-N16	-3.11	102.70	111.16
3	A	401	SNA	C47-C48-C45	2.85	122.65	118.17
3	A	401	SNA	O38-P35-O36	-2.76	106.86	113.06
3	A	401	SNA	C14-C15-N16	2.73	116.54	110.79
3	A	401	SNA	P1-C5-C8	2.71	117.10	108.95
3	A	401	SNA	C46-C48-C47	-2.62	117.14	120.89
3	A	401	SNA	C15-C17-N18	-2.62	111.26	116.54
3	A	401	SNA	F7-C5-F6	2.58	109.99	106.73
3	A	401	SNA	O37-P35-O36	-2.54	107.36	113.06
3	A	401	SNA	F40-C39-C43	-2.45	107.33	110.49
3	A	401	SNA	C48-C46-C33	-2.07	106.44	112.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	SNA	C42-C43-C39	2.04	121.69	119.84

There are no chirality outliers.

All (7) torsion outliers are listed below:

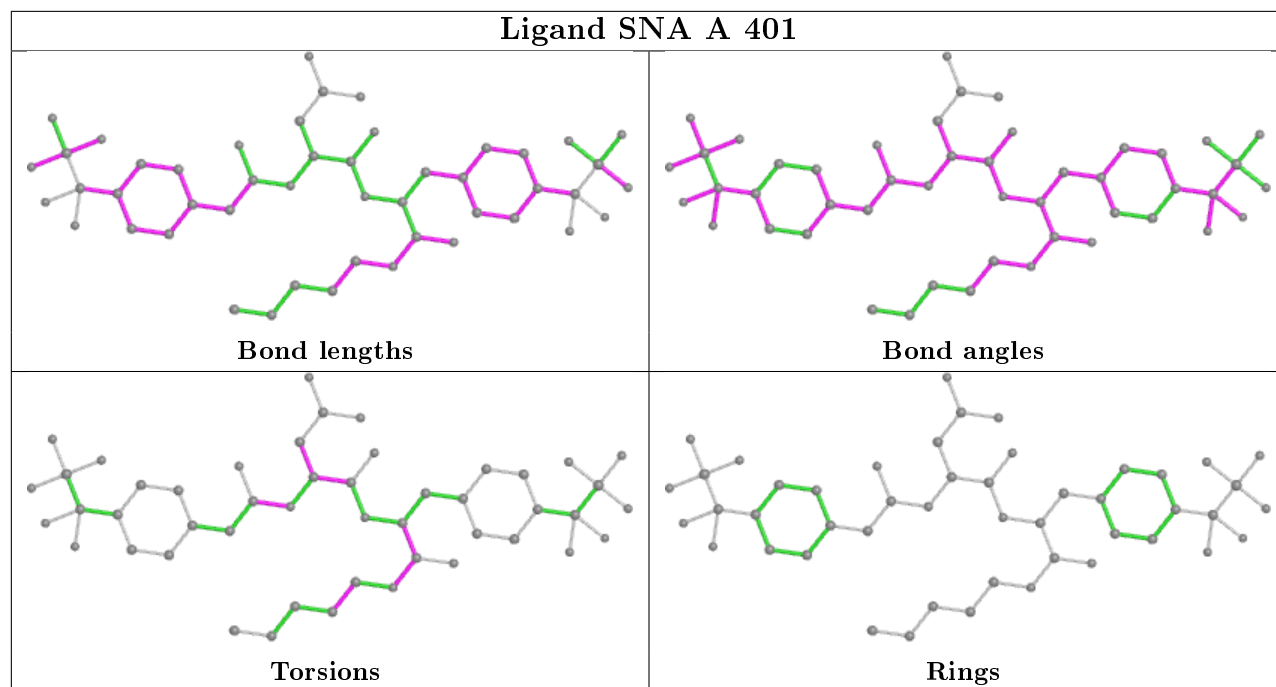
Mol	Chain	Res	Type	Atoms
3	A	401	SNA	C15-C17-N18-C20
3	A	401	SNA	N25-C26-C27-C28
3	A	401	SNA	O34-C33-N25-C26
3	A	401	SNA	N18-C20-C21-C22
3	A	401	SNA	C27-C26-C31-N16
3	A	401	SNA	C27-C26-C31-O32
3	A	401	SNA	N16-C15-C17-N18

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	SNA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.